

ER PROGRAM DATA ASSESSMENT
SUMMARY REPORT FORM

Batch No. 8910L081 Site Solar Ponds
Laboratory Roy F. Weston - Lionville No. of Samples/Matrix 4/Water
SOW # 10/86 (Rev. 2/88) Reviewer Org. TechLaw, Inc.
Sample Numbers SW094007, TB101089007, SW095007, SW093007

Data Assessment Summary

	VOA	Comments
1. Holding Times	<u>A</u>	<u>Action Items 1,2; Comment 1</u>
2. GC/MS Tune/Instr. Perf.	<u>V</u>	
3. Calibrations	<u>A</u>	<u>Action Items 3,4; Comment 2</u>
4. Blanks	<u>A</u>	<u>Action Item 5</u>
5. Surrogates	<u>A</u>	<u>Action Items 6,7</u>
6. Matrix Spike/Dup.	<u>X</u>	<u>Comment 3</u>
7. Other QC	<u>V</u>	
8. Internal Standards	<u>A</u>	<u>Action Items 6,7</u>
9. Compound Identification	<u>X</u>	<u>Action Items 6,8; Comments 4,5</u>
10. System Performance	<u>X</u>	<u>Comment 6</u>
11. Overall Assessment	<u>A</u>	<u>Data acceptable with qualifications.</u>

V = Data had no problems.

A = Data acceptable but qualified due to problems.

R = Data rejected.

X = Problems, but do not affect data.

Data Quality: Data contained in this batch were reviewed and found to be acceptable with qualifications. Acceptable,
qualified data may be used provided that individual values impacted by the "Action Items" listed below are appropriately flagged.
(Refer to attached Results Summary Table.)

REVIEWED FOR CLASSIFICATION
By [Signature]
Date 10/25/91

Action Items: 1) Non-detected results for aromatic compounds in samples SW093007, SW094007, and TB101089007 are estimated and undetected (UJ) because holding times exceeded seven days.

2) All results in the dilution of sample SW095007 are rejected (R) because the holding time exceeded fourteen days and the duration from VTSR to sample analysis exceeded ten days.

3) 4-Methyl-2-pentanone's RRF were less than 0.05 in all calibrations and 2-Butanone had either RRFs less than 0.05, a %RSD or %D greater than 50% in the initial and continuing calibrations. As a result, the non-detected results for these two compounds in all analyses are rejected (R).

4) Acetone and Methylene Chloride had %RSDs exceeding 50% in the 10/18/89 initial calibration and Acetone's %D exceeded 50% in the 10/20/89 continuing calibration. Therefore, the non-detected results for Methylene Chloride in sample SW094007 and Acetone in sample SW095007 are rejected (R). The positive results for Acetone in the other three samples are estimated (J). Positive results for Methylene Chloride in samples SW093007 and TB101089007 would be estimated (J) had blank criteria been met. See Action Item 5. The positive result for Methylene Chloride in sample SW095007 would be estimated however, all results are rejected (R) for this sample. See Action Item 6.

5) As a result of method blank contamination, the positive results for Methylene Chloride in samples SW093007 and TB101089007 are estimated and undetected (UJ) according to the Functional Guidelines criteria (5x and 10x rules).

6) Surrogate recoveries exceeded 300% and all three internal standard areas exceeded criteria in sample SW095007. In addition, unknown peaks were evident in the chromatogram, including one enormous peak, all of which were not quantitated or reported. Furthermore, the quantitation report had positive results for several compounds that were changed to non-detects without being initialed or dated. None of these peaks or compounds, including Methylene Chloride which exceeded calibration range, were evident in the dilution. As a result of these contributory problems, all results in sample SW095007 are rejected (R).

7) All three internal standard areas exceeded criteria in samples SW093007, SW094007, and TB101089007 and sample SW094007 also had surrogate recoveries outside criteria. Therefore, all results in these three samples are estimated (J) or estimated and undetected (UJ).

Action Items: (cont) 8) The positive result for Chloroform in sample SW094007 is estimated (J) because mass spectral data was not submitted to confirm its identity. The value is reported because of a good RRT match with the 12 hour standard.

Comments: 1) Sample reanalyses were conducted beyond the fourteen day and ten day contract holding times. Therefore, the original analyses data are reported on the Data Summary Table.

2) Other compounds whose %RSD or %D exceeded criteria in the calibrations were undetected in the samples and no action is necessary.

3) No MS/MSD analysis was performed with these samples to determine if possible matrix effects contributed to surrogate and internal standard problems. No further action is necessary as all sample results are qualified.

4) The reference spectrum for Acetone did not match NBS standard spectrum.

5) Vinyl Acetate's identity in sample SW095007 is not confirmed by the mass spectrum provided. However all results in this sample are rejected.

6) The surrogate and internal standard area difficulties may indicate a serious system problem.

Note: Data Summary Tables are attached.

William T Fee
Reviewer Signature

6/25/80
Date

SITE NAME: Solar Ponds

CLP VOLATILE ORGANIC ANALYSIS: Low Water

ANALYTICAL RESULTS (ppb)

Sample Location	Sample Number	SW093007	SW094007	TB101088007	SW095007	VBLK058	SW095007DL
Sampling Date	10/10/89	10/10/89	10/10/89	10/10/89	10/10/89	Method Blank	2x Dilution
Remarks	Method Blank			Trip Blank			
Volatiles	CRQL	DQ	DQ	DQ	DQ	DQ	DQ
Compound	ug/L (ppb)						
Chloromethane	10	10 U A	10 U A	10 U A	10 U R		20 U R
Bromomethane	10	10 U A	10 U A	10 U A	10 U R		20 U R
Vinyl chloride	10	10 U A	10 U A	10 U A	10 U R		20 U R
Chloroethane	10	10 U A	10 U A	10 U A	10 U R		20 U R
Methylene chloride	5	6 ppb	5 U R	5 U A	240 R	3 ppb	10 U R
Acetone	10	24 J A	7 J A	7 J A	10 U R		20 U R
Carbon disulfide	5	5 U A	5 U A	5 U A	5 U R		10 U R
1,1-Dichloroethene	5	5 U A	5 U A	5 U A	5 U R		10 U R
1,1-Dichloroethane	5	5 U A	5 U A	5 U A	5 U R		10 U R
1,2-Dichloroethene (Total)	5	5 U A	5 U A	5 U A	5 U R		10 U R
Chloroform	5	5 U A	1 J A	5 U A	2 R		10 U R
1,2-Dichloroethane	5	5 U A	5 U A	5 U A	5 U R		10 U R
2-Butanone	10	10 U R	10 U R	10 U R	10 U R		20 U R
1,1,1-Trichloroethane	5	5 U A	5 U A	5 U A	5 U R		10 U R
Carbon tetrachloride	5	5 U A	5 U A	5 U A	5 U R		10 U R
Vinyl acetate	10	10 U A	10 U A	10 U A	2 R		20 U R
Bromodichloromethane	5	5 U A	5 U A	5 U A	5 U R		10 U R
1,2-Dichloropropane	5	5 U A	5 U A	5 U A	5 U R		10 U R
cis-1,3-Dichloropropene	5	5 U A	5 U A	5 U A	5 U R		10 U R
Trichloroethene	5	5 U A	5 U A	5 U A	3 R		10 U R
Dibromodichloromethane	5	5 U A	5 U A	5 U A	5 U R		10 U R
1,1,1,2-Trichloroethane	5	5 U A	5 U A	5 U A	5 U R		10 U R
Benzene	5	5 U A	5 U A	5 U A	5 U R		10 U R
trans-1,3-Dichloropropene	5	5 U A	5 U A	5 U A	5 U R		10 U R
Bromoforn	5	5 U A	5 U A	5 U A	5 U R		10 U R
4-Methyl-2-pentanone	10	10 U R	10 U R	10 U R	10 U R		20 U R
2-Hexanone	10	10 U A	10 U A	10 U A	10 U R		20 U R
Tetrachloroethene	5	5 U A	5 U A	5 U A	5 U R		10 U R
1,1,2,2-Tetrachloroethane	5	5 U A	5 U A	5 U A	5 U R		10 U R
Toluene	5	5 U A	5 U A	5 U A	5 U R		10 U R
Chlorobenzene	5	5 U A	5 U A	5 U A	5 U R		10 U R
Ethylbenzene	5	5 U A	5 U A	5 U A	5 U R		10 U R
Styrene	5	5 U A	5 U A	5 U A	5 U R		10 U R
Xylenes (Total)	5	5 U A	5 U A	5 U A	5 U R		10 U R
Total Organic Concentration (ppb)	6	24	8	7	247	3	0

U Indicates the compound was not detected above the Required Quantitation Limit.

J Quantitation is approximate due to limitations identified during the quality control review.

E Exceeds calibration range, dilute & reanalyze.

CRQL Contract Required Quantitation Limit in Micrograms per Liter (ug/L), Parts per billion (ppb).

DQ Data Qualifier

V Valid

A Acceptable with qualifications

R Rejected

ER DEPARTMENT DATA ASSESSMENT
SUMMARY REPORT FORM

Batch No. 8910L081 Site Solar Ponds
Laboratory Roy F. Weston - Lionville No. of Samples/Matrix 3/Water
SOW # 10/86 (Rev. 2/88) Reviewer Org. TechLaw, Inc.
Sample Numbers SW094007, SW095007, SW093007

Data Assessment Summary

	BNA	Comments
1. Holding Times	<u>V</u>	<u></u>
2. GC/MS Tune/Instr. Perf.	<u>V</u>	<u></u>
3. Calibrations	<u>A</u>	<u>Action Item 1; Comment 1</u>
4. Blanks	<u>A</u>	<u>Action Item 2; Comment 2</u>
5. Surrogates	<u>X</u>	<u>Comment 3</u>
6. Matrix Spike/Dup.	<u>X</u>	<u>Comment 4</u>
7. Other QC	<u>V</u>	<u></u>
8. Internal Standards	<u>V</u>	<u></u>
9. Compound Identification	<u>X</u>	<u>Comment 5</u>
10. System Performance	<u>V</u>	<u></u>
11. Overall Assessment	<u>A</u>	<u>Data Acceptable with Qualifications</u>

V = Data had no problems.

A = Data acceptable but qualified due to problems.

R = Data rejected.

X = Problems, but do not affect data.

Data Quality: Data contained in this batch were reviewed and found to be acceptable with qualifications. Acceptable, qualified data may be used provided that individual values impacted by the "Action Items" listed below are appropriately flagged.
(Refer to attached Data Summary Tables.)

Action Items: 1) The %RSD for Benzyl Alcohol exceeded 50% in the initial calibration. The non-detected Benzyl Alcohol results are rejected (R) in all samples.

2) As a result of method blank contamination, the positive results for Di-n-butylphthalate and Bis(2-ethylhexyl)phthalate in all samples is estimated and undetected (UJ) according to the Functional Guidelines criteria.

Comments: 1) Several compounds exceeded the %RSD and %D criteria in the 10/17/89 initial and the 10/19/89 continuing calibrations. Furthermore, Benzoic Acid's RRF was less than 0.05 in this continuing calibration. The only sample affected is the method blank and, therefore, data is not qualified.

2) The method blank was analyzed on a different instrument than the samples. Contamination found in the blank only relates to possible contaminants in the samples introduced in the extraction process and not the analysis itself.

3) Two surrogates fell outside the % recovery criteria for the blank matrix spike duplicate sample analysis.

4) A blank should not be used for the MS/MSD analysis since it is not representative of the sample matrix. Moreover, six compounds exceeded the % recovery in the BS and BSD samples. No action is necessary because results are not qualified based on MS/MSD data alone.

5) Samples SW093007 and SW095007 contained four TICs and sample SW094007 contained five TICs.

Note: Data Summary Tables are attached.

Lisa Contreras - Hendler
Reviewer Signature

7-11-90
Date

SITE NAME: Solar Ponds

CLP SEMIVOLATILE ANALYSIS: Low Water

ANALYTICAL RESULTS (ug/L)

Sample Location	Sample Number	SBLK 967	SW034007	SW095007	SW083007			
Sample Number								
Sampling Date			10/10/89	10/10/89	10/10/89			
Remarks		Method Blank						
Semivolatiles (BVA)	CRQL	DQ	DQ	DQ	DQ			
Organic Compound	ug/L							
Phenol	10		10 U V	10 U V	10 U V			
Bis(2 - Chloroethyl)ether	10		10 U V	10 U V	10 U V			
2 - Chlorophenol	10		10 U V	10 U V	10 U V			
1,3 - Dichlorobenzene	10		10 U V	10 U V	10 U V			
1,4 - Dichlorobenzene	10		10 U V	10 U V	10 U V			
Benzyl alcohol	10		10 U R	10 U R	10 U R			
1,2 - Dichlorobenzene	10		10 U V	10 U V	10 U V			
2 - Methylphenol	10		10 U V	10 U V	10 U V			
Bis(2-chloroisopropyl)ether	10		10 U V	10 U V	10 U V			
4 - Methylphenol	10		10 U V	10 U V	10 U V			
N-Nitroso-d-n-propylamine	10		10 U V	10 U V	10 U V			
Hexachloroethane	10		10 U V	10 U V	10 U V			
Nitrobenzene	10		10 U V	10 U V	10 U V			
Isophorone	10		10 U V	10 U V	10 U V			
2 - Nitrophenol	10		10 U V	10 U V	10 U V			
2,4 - Dimethylphenol	10		10 U V	10 U V	10 U V			
Benzoic Acid	50		50 U V	50 U V	50 U V			
Bis(2-Chloroethoxy)methane	10		10 U V	10 U V	10 U V			
2,4 - Dichlorophenol	10		10 U V	10 U V	10 U V			
1,2,4 - Trichlorobenzene	10		10 U V	10 U V	10 U V			
Naphthalene	10		10 U V	10 U V	10 U V			
4 - Chloroaniline	10		10 U V	10 U V	10 U V			
Hexachlorobutadiene	10		10 U V	10 U V	10 U V			
4-Chloro-3-methylphenol	10		10 U V	10 U V	10 U V			
2-Methylnaphthalene	10		10 U V	10 U V	10 U V			
Hexachlorocyclopentadiene	10		10 U V	10 U V	10 U V			
2,4,6-Trichlorophenol	10		10 U V	10 U V	10 U V			
2,4,5-Trichlorophenol	50		50 U V	50 U V	50 U V			
2-Chloronaphthalene	10		10 U V	10 U V	10 U V			
2-Nitroaniline	50		50 U V	50 U V	50 U V			
Dimethyl phthalate	10		10 U V	10 U V	10 U V			
Acenaphthylene	10		10 U V	10 U V	10 U V			
2,6-Dinitrotoluene	10		10 U V	10 U V	10 U V			
3-Nitroaniline	50		50 U V	50 U V	50 U V			

E Exceeds calibration range.

U Indicates the compound was not detected above the Instrument Quantitation Limit.

J Quantitation is approximate due to limitations identified during the quality control review.

CRQL Contract Required Quantitation Limit in Micrograms per Liter (ug/L), Parts per Billion (ppb).

DQ Data Qualifier

V Valid

A Acceptable with qualifications

R Rejected

OL081.WK1

SITE NAME: Solar Ponds

CLP SEMIVOLATILE ANALYSIS: Low Water

ANALYTICAL RESULTS (ug/L)

Sample Location	SBLK 967	SW094007	SW095007	SW093007		
Sample Number		10/10/89	10/10/89	10/10/89		
Sampling Date						
Remarks	Method Blank					
Semivolatiles (BNA)						
Organic Compound	CRQL ug/L	DQ	DQ	DQ		
Acenaphthene	10	10 U V	10 U V	10 U V	10 U V	DQ
2,4-Dinitrophenol	50	50 U V	50 U V	50 U V	50 U V	
4-Nitrophenol	50	50 U V	50 U V	50 U V	50 U V	
Dibenzofuran	10	10 U V	10 U V	10 U V	10 U V	
2,4-Dinitrotoluene	10	10 U V	10 U V	10 U V	10 U V	
Diethyl phthalate	10	10 U V	10 U V	10 U V	10 U V	
4-Chlorophenyl-phenyl ether	10	10 U V	10 U V	10 U V	10 U V	
Fluorene	10	10 U V	10 U V	10 U V	10 U V	
4-Nitroaniline	50	50 U V	50 U V	50 U V	50 U V	
4,6-Dinitro-2-methylphenol	50	50 U V	50 U V	50 U V	50 U V	
N-Nitrosodiphenylamine	10	10 U V	10 U V	10 U V	10 U V	
4-Bromophenyl phenyl ether	10	10 U V	10 U V	10 U V	10 U V	
Hexachlorobenzene	10	10 U V	10 U V	10 U V	10 U V	
Pentachlorophenol	50	50 U V	50 U V	50 U V	50 U V	
Phenanthrene	10	10 U V	10 U V	10 U V	10 U V	
Anthracene	10	10 U V	10 U V	10 U V	10 U V	
Di-n-butyl phthalate	10	10 U V	10 U V	10 U V	10 U V	
Fluoranthene	10	10 U V	10 U V	10 U V	10 U V	
Pyrene	10	10 U V	10 U V	10 U V	10 U V	
Butylbenzyl phthalate	10	10 U V	10 U V	10 U V	10 U V	
3,3'-Dichlorobenzidine	20	20 U V	20 U V	20 U V	20 U V	
Benzo(a)anthracene	10	10 U V	10 U V	10 U V	10 U V	
Chrysene	10	10 U V	10 U V	10 U V	10 U V	
Bis(2-ethylhexyl)phthalate	10	14 U A	24 U A	15 U A	15 U A	
Di-n-octyl phthalate	10	10 U V	1 J A	10 U V	10 U V	
Benzo(b)fluoranthene	10	10 U V	10 U V	10 U V	10 U V	
Benzo(k)fluoranthene	10	10 U V	10 U V	10 U V	10 U V	
Benzo(a)pyrene	10	10 U V	10 U V	10 U V	10 U V	
Indeno(1,2,3-cd)pyrene	10	10 U V	10 U V	10 U V	10 U V	
Dibenz(a,h)anthracene	10	10 U V	10 U V	10 U V	10 U V	
Benzo(g,h,i)perylene	10	10 U V	10 U V	10 U V	10 U V	
Total semivolatile BNA concentration (ppb)	6	0	1	0	0	

E Exceeds calibration range.

U Indicates the compound was not detected above the Instrument Quantitation Limit.

J Quantitation is approximate due to limitations identified during the quality control review.

CRQL Contract Required Quantitation Limit in Micrograms per Liter (ug/L), Parts per Billion (ppb).

DQ
V
A
R

Data Qualifier
Valid
Acceptable with qualifications
Rejected

ER DEPARTMENT DATA ASSESSMENT
SUMMARY REPORT FORM

Batch No. 8910L081 Site Solar Ponds
Laboratory Roy F. Weston - Lionville No. of Samples/Matrix 3/Water
SOW # 10/86 (Rev. 2/88) Reviewer Org. TechLaw, Inc.
Sample Numbers SW094007, SW095007, SW093007

Data Assessment Summary

	Pesticides/PCB	Comments
1. Holding Times	<u>V</u>	<u></u>
2. Instrument Performance	<u>X</u>	<u>Comment 1</u>
3. Calibrations	<u>X</u>	<u>Comment 2</u>
4. Blanks	<u>V</u>	<u></u>
5. Surrogates	<u>X</u>	<u>Comment 3</u>
6. Matrix Spike/Dup.	<u>X</u>	<u>Comment 4</u>
7. Other QC	<u>X</u>	<u>Comment 5</u>
8. Compound Identification	<u>X</u>	<u>Comment 6</u>
9. System Performance	<u>V</u>	<u></u>
10. Overall Assessment	<u>V</u>	<u>Data is valid.</u>

V = Data had no problems.

A = Data acceptable but qualified due to problems.

R = Data rejected.

X = Problems, but do not affect data.

Data Quality: Data contained in this batch were reviewed and found to be valid. (Refer to attached Data Summary Table.)

Comments: 1) In the continuing calibration of 10/26/89 (0419) three compounds were outside of their retention time windows. In the continuing calibration of 10/26/89 (0452) one compound was outside of its retention time window. These calibrations were on the confirmation column. Expanded windows were used to evaluate the chromatograms and no action is necessary.

2) The calibration factors of the multiresponse compounds (Aroclors and Toxaphene) are not reproducible. It is suspected that multiple peaks were used to calculate the calibration factors. There were no positive results for these compounds and no action is necessary.

3) Surrogate Recoveries were high in the method blank PBLK966 and sample SW093007. No action is necessary.

4) No matrix spike/matrix spike duplicate data was provided with this batch. Although blank spike data was included, this does not represent a sample matrix. In this blank spike data, Lindane's (gamma-BHC) RPD and Percent Recoveries in both the BS/BSD exceeded the criteria. No action is necessary because results are not qualified due to MS/MSD alone.

5) Although 990 milliliters of sample were extracted, the CROLs were not adjusted accordingly. Therefore, the CROLs were recalculated and the changed results reported on the Data Summary Table.

6) The positive result for DDT in the method blank is questioned. The retention time of DDT on the confirmation column shifted, which is not a problem; however, it shifted early rather than late like the surrogate. In addition the quantities generated from the two columns were very different. Nevertheless, the result is well below the CROL and no action is taken.

Note: Data Summary Tables are attached.

Jill Gascher
Reviewer Signature

7-10-90
Date

SITE NAME: Solar Ponds

CLP PESTICIDES/PCB ANALYSIS: Low Water

ANALYTICAL RESULTS (ug/L)

Sample Location	PBLK966	SW094007	SW095007	SW093007				
Sample Number		10/10/89	10/10/89	10/10/89				
Sample Date								
Remarks	Method Blank							
Chlorinated Pesticides	CRQL ug/L	DQ	DQ	DQ	DQ			
alpha - BHC	0.050	0.051 U V	0.051 U V	0.051 U V	0.051 U V			
beta - BHC	0.050	0.051 U V	0.051 U V	0.051 U V	0.051 U V			
delta - BHC	0.050	0.051 U V	0.051 U V	0.051 U V	0.051 U V			
gamma - BHC (Lindane)	0.050	0.051 U V	0.051 U V	0.051 U V	0.051 U V			
Heptachlor	0.050	0.051 U V	0.051 U V	0.051 U V	0.051 U V			
Aldrin	0.050	0.051 U V	0.051 U V	0.051 U V	0.051 U V			
Heptachlor epoxide	0.050	0.051 U V	0.051 U V	0.051 U V	0.051 U V			
Endosulfan I	0.050	0.051 U V	0.051 U V	0.051 U V	0.051 U V			
Dieldrin	0.10	0.10 U V	0.10 U V	0.10 U V	0.10 U V			
4,4' - DDE	0.10	0.10 U V	0.10 U V	0.10 U V	0.10 U V			
Endrin	0.10	0.10 U V	0.10 U V	0.10 U V	0.10 U V			
Endosulfan II	0.10	0.10 U V	0.10 U V	0.10 U V	0.10 U V			
4,4' - DDD	0.10	0.10 U V	0.10 U V	0.10 U V	0.10 U V			
Endosulfan sulfate	0.10	0.10 U V	0.10 U V	0.10 U V	0.10 U V			
4,4' - DDT	0.10	0.10 U V	0.10 U V	0.10 U V	0.10 U V			
Methoxychlor	0.50	0.51 U V	0.51 U V	0.51 U V	0.51 U V			
Endrin Ketone	0.10	0.10 U V	0.10 U V	0.10 U V	0.10 U V			
alpha - Chlordane	0.50	0.51 U V	0.51 U V	0.51 U V	0.51 U V			
gamma - Chlordane	0.50	0.51 U V	0.51 U V	0.51 U V	0.51 U V			
Toxaphene	1.0	1.0 U V	1.0 U V	1.0 U V	1.0 U V			
Aroclor - 1016	0.50	0.51 U V	0.51 U V	0.51 U V	0.51 U V			
Aroclor - 1221	0.50	0.51 U V	0.51 U V	0.51 U V	0.51 U V			
Aroclor - 1232	0.50	0.51 U V	0.51 U V	0.51 U V	0.51 U V			
Aroclor - 1242	0.50	0.51 U V	0.51 U V	0.51 U V	0.51 U V			
Aroclor - 1248	0.50	0.51 U V	0.51 U V	0.51 U V	0.51 U V			
Aroclor - 1254	1.0	1.0 U V	1.0 U V	1.0 U V	1.0 U V			
Aroclor - 1260	1.0	1.0 U V	1.0 U V	1.0 U V	1.0 U V			
Total Chlorinated Pesticides (ppb)	0.80	0	0	0	0			

E Exceeds calibration range.

U Indicates the compound was not detected above the Instrument Quantitation Limit

J Quantitation is approximate due to limitations identified during the quality control review

CRQL Contract Required Detection Limit in Micrograms per Liter (ug/L), Parts per Billion (ppb)

DQ Data Qualifier

V Valid

A Acceptable with qualifications

R Rejected

10L081/pcb

ER PROGRAM DATA ASSESSMENT
SUMMARY REPORT FORM

Batch No. 8910L081 Site Area 6 - Solar Ponds
Laboratory Roy F. Weston - Lionville No. of Samples/Matrix 6/Water
SOW # 7/87 Reviewer Org. TechLaw, Inc.
Sample Numbers SW094007 (total), SW095007 (total), SW093007 (total) SW094007 (soluble),
SW095007 (soluble), SW093007 (soluble)

Data Assessment Summary

	ICP	AA	Hg	CN	Comments
1. Holding Times	<u>V</u>	<u>V</u>	<u>V</u>	<u>V</u>	
2. Calibrations	<u>A</u>	<u>V</u>	<u>V</u>	<u>V</u>	<u>Action Items 1-4</u>
3. Blanks	<u>A</u>	<u>A</u>	<u>V</u>	<u>V</u>	<u>Action Items 5-10</u>
4. ICP Interference Check Sample	<u>A</u>	<u>N/A</u>	<u>N/A</u>	<u>N/A</u>	<u>Action Items 11-13</u>
5. Lab Control Sample Results	<u>V</u>	<u>V</u>	<u>V</u>	<u>V</u>	
6. Duplicate Sample Results	<u>V</u>	<u>V</u>	<u>V</u>	<u>V</u>	
7. Matrix Spike Sample Results	<u>V</u>	<u>A</u>	<u>V</u>	<u>V</u>	<u>Action Items 14-17</u>
8. Method of Standard Addition	<u>N/A</u>	<u>V</u>	<u>N/A</u>	<u>N/A</u>	
9. Serial Dilution	<u>A</u>	<u>N/A</u>	<u>N/A</u>	<u>N/A</u>	<u>Action Item 18</u>
10. Sample Verification	<u>V</u>	<u>V</u>	<u>V</u>	<u>V</u>	
11. Other QC	<u>V</u>	<u>V</u>	<u>V</u>	<u>V</u>	
12. Overall Assessment	<u>A</u>	<u>A</u>	<u>V</u>	<u>V</u>	<u>Data valid, or acceptable with qualifications</u>

V = Data had no problems.

A = Data acceptable but qualified due to problems.

R = Data rejected.

X = Problems, but do not affect data.

N/A = Not applicable.

Data Quality: Data contained in this batch were reviewed and found to be valid, or acceptable with qualifications. Acceptable,
qualified data may be used provided that individual values impacted by the "Action Items" listed below are appropriately flagged.
(Refer to attached Results Summary Tables).

Action Items: 1) All non-detect Molybdenum values are estimated and undetected (UJ) because the CRI recovery criteria were not met.

2) The Zinc value for SW093007 (soluble) is estimated (J) because the CRI recovery criteria were not met.

3) The non-detect Lithium aqueous values for SW093007 (total and soluble) are estimated and undetected (UJ) because the CRI recovery criteria were not met.

4) All non-detect Tin values except SW093007 (total) are estimated and undetected (UJ) and the value for SW093007 (total) is estimated (J) because the CRI recovery criteria were not met.

5) All Aluminum values are estimated and undetected (UJ) because Aluminum values >IDL were found in the blanks.

6) The Cobalt values for SW094007 (total) and SW093007 (soluble) are estimated and undetected (UJ) because Cobalt values >IDL were found in the blanks.

7) All Copper values are estimated and undetected (UJ) because Copper values >IDL were found in the blanks.

8) The Potassium value for SW093007 (total and soluble) is estimated and undetected (UJ) because Potassium values >IDL were found in the blanks.

9) The Antimony values for SW094007 (total and soluble), SW095007 (total), and SW093007 (total) are estimated and undetected (UJ) because Antimony values >IDL were found in the blanks.

10) The Thallium value for SW095007 (soluble) is rejected (R) because of negative bias indicated in the blanks.

11) The Silver values for SW094007 (total and soluble) and SW095007 (total and soluble) are rejected (R) because the interference check sample result was outside control limits.

12) The Chromium values for SW094007 (total and soluble) and SW095007 (total and soluble) are rejected (R) because the interference check sample result was outside control limits.

13) The Manganese, Beryllium, and Zinc values for SW094007 (total and soluble) and SW095007 (total and soluble) are estimated (J) because the interference check sample result was outside control limits.

Action Items: (cont) 14) All Lead values except SW093007 (soluble) are estimated and undetected (UJ)
and the value for SW093007 (soluble) is estimated (J) because the pre-digestion matrix spike recovery criteria were
not met.

15) All Selenium values are estimated (J) because the pre-digestion matrix spike recovery criteria were not
met.

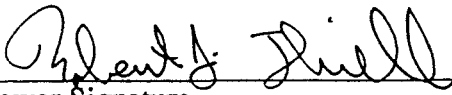
16) All Thallium values except SW093007 (total) are estimated and undetected (UJ) because the pre-digestion
matrix spike recovery criteria were not met.

17) The non-detect Arsenic values for SW094007 (soluble), SW095007 (soluble), and SW093007 (soluble) are
estimated and undetected (UJ) because the post-digestion matrix spike recovery criteria were not met.

18) All Calcium and Magnesium values are estimated (J) because the ICP serial dilution recovery criteria were
not met.

Comments: None

Note: Data Summary Tables are attached.


Reviewer Signature

2/21/90
Date

SITE NAME: Site Background Characterization

CLP WATER INORGANIC ANALYSIS: Low Water

ANALYTICAL RESULTS (ug/L)

Sample Location	SW093007	SW093007	SW094007	SW094007	SW095007	SW095007							
Sample Number	10/1089	10/1089	10/1089	10/1089	10/1089	10/1089							
Sample Date	Total	Soluble	Total	Soluble	Total	Soluble							
Remarks													
Inorganic Analyte	DQ	DQ	DQ	DQ	DQ	DQ							
DL ug/L													
Aluminum	126 UJ	A	164 UJ	A	267 UJ	A	260 UJ	A	277 UJ	A	293 UJ	A	
Antimony	26.2 UJ	A	22.0 U	V	42.0 UJ	A	28.4 UJ	A	39.9 UJ	A	22.0 U	V	
Arsenic	2.0 U	V	2.0 UJ	A	2.0 U	V	2.0 UJ	A	2.0 U	V	2.0 UJ	A	
Barium	137	V	134	V	177	V	170	V	176	V	177	V	
Beryllium	1.0 U	V	1.0 U	V	1.3 J	A	1.3 J	A	1.3 J	A	1.3 J	A	
Cadmium	3.0 U	V	3.0 U	V	3.0 U	V	3.0 U	V	3.0 U	V	3.0 U	V	
Calcium	112000 J	A	109000 J	A	389000 J	A	373000 J	A	392000 J	A	396000 J	A	
Cesium	2500 U	V	2500 U	V	2500 U	V	2500 U	V	2500 U	V	2500 U	V	
Chromium	2.0 U	V	2.0 U	V	2.0 U	R	2.0 U	R	2.0 U	R	2.0 U	R	
Cobalt	4.0 U	V	4.4 UJ	A	5.1 UJ	A	4.0 U	V	4.0 U	V	4.0 U	V	
Copper	11.5 UJ	A	16.1 UJ	A	24.3 UJ	A	20.7 UJ	A	25.1 UJ	A	23.2 UJ	A	
Iron	551	V	1220	V	104	V	104	V	115	V	108	V	
Lead	3.0 UJ	A	85.6 J	A	3.0 UJ	A	3.0 UJ	A	3.0 UJ	A	3.0 UJ	A	
Lithium	100 UJ	A	100 UJ	A	457	V	434	V	461	V	464	V	
Magnesium	30400 J	A	29200 J	A	106000 J	A	102000 J	A	106000 J	A	107000 J	A	
Manganese	15	870	V	847	V	18.1 J	A	14.8 J	A	18.5 J	A	18.1 J	A
Mercury	0.2	0.20 U	V	0.20 U	V	0.20 U	V	0.20 U	V	0.20 U	V	0.20 U	V
Molybdenum	200	100 UJ	A	100 UJ	A	100 UJ	A	100 UJ	A	100 UJ	A	100 UJ	A
Nickel	40	7.0 U	V	7.0 U	V	9.3	V	7.0 U	V	10.3	V	7.0 U	V
Potassium	5000	3640 UJ	A	4300 UJ	A	91000	V	86700	V	91300	V	91200	V
Selenium	5	37.4 J	A	62.2 J	A	16.4 J	A	15.8 J	A	15.0 J	A	9.2 J	A
Silver	10	3.0 U	V	3.0 U	V	3.0 U	R	3.0 U	R	3.0 U	R	3.0 U	R
Sodium	5000	61900	V	59600	V	563000	V	571000	V	590000	V	578000	V
Strontium	200	779	V	745	V	3120	V	3000	V	3120	V	3140	V
Thallium	10	4.0 UJ	A	4.0 UJ	A	40.0 UJ	A	4.0 UJ	A	4.0 UJ	A	4.0 U	R
Tin	200	100 UJ	A	100 UJ	A	136 J	A	114 J	A	128 J	A	131 J	A
Vanadium	50	5.0 U	V	5.0 U	V	5.0 U	V	5.0 U	V	5.0 U	V	5.0 U	V
Zinc	20	65.6	V	54.3 J	A	18.7 J	A	23.4 J	A	23.4 J	A	24.8 J	A
Cyanide	10	N/R	N/R	N/R	10.0 U	V	N/R	N/R	10.0 U	V	N/R	N/R	N/R

E Estimated by the Laboratory

U Indicates the compound was not detected above the Instrument Quantitation Limit

J Quantitation is approximate due to limitations identified during the quality control review

DL Detection Limit in Micrograms per Liter (ug/L)

N/R Not reported

DQ Data Qualifier

V Valid

A Acceptable with qualifications

R Rejected

L0811Ueg04]

Lot 101089001

Custody Transfer Record/Lab Work Request

WESTON Analytics Use Only
8910L081

Client Backwell
Work Order 2029-3304
Date Rec'd. 10/12/89 Date Due 11/1/89
RFW Contact M. Carbaugh
Client Contact/Phone _____



Client: <u>Rockwell</u>		Work Order: <u>2029-3304</u>		Date Rec'd: <u>10/2/89</u>		Date Due: <u>11/09/89</u>	
RFW Contact: <u>M. Carthage</u>		Client Contact/Phone: _____					

WA Use Only	Lab ID	Client ID/Description	Matrix	Date Collected	Analyses Requested	Refrigerator #	#/Type Container	Volume	Preservative
	001X	SW094007 radioactive	W	10-10-89					
	002X	TB101089007	W						
	003X	SW095007 radioactive	W						
	004X	SW093007 radioactive	W						
	005	SW094007 radioactive							
		TB101089007-KS							
	006	SW095007 radioactive							
	007	SW093007 radioactive							

Item/Reason	Relinquished by	Received by	Date	Time
ALL	James P. [Signature]	James P. [Signature]	11/18/89	1420
ALL	James P. [Signature]	James P. [Signature]	10/18/89	1420
ALL	James P. [Signature]	James P. [Signature]	10/12/89	9:30 AM
ALL	James P. [Signature]	James P. [Signature]	11/05/89	0830
ALL	James P. [Signature]	James P. [Signature]	11/05/89	1650

Matrix:	W - Water	DS - Drum Solids	X - Other
S - Soil	O - Oil	DL - Drum Liquids	
SE - Sediment	A - Air	F - Fish	
SO - Solid	WI - Wipe	L - EP/TCLP Leachate	

Special Instructions: 1 - vos 2 - bna 3 - pest/pcb 4 - cyanide
 5 - filtered tcl metals Mo, Sr, Ca, Li, Sn
 6 - unfiltered tcl metals Mo, Sr, Ca, Li, Sn

Retention 10/20/89 Sample
 #48 Tape 10/20/89 Radiactive
 previous Scott.

WESTON ANALYTICS
TOTAL RADIOCHEMISTRY
DATA SUMMARY REPORT

Client: ROCKWELL (ROCKY FLATS)

Page: 3

Sample Information

	8910505	89106787(?)	89106081
RFW Batch ID:	9003-798-1	9003-798-2	9003-803-5
Customer ID:	GWSF06891089001	604861089004	SH094007
Collection Date:	10/05/89	10/06/89	10/10/89
Matrix:	Water	Water	Water

Radio Chemistry

Gross Alpha.....	1.3 ± 0.7	pci/l 1.0	6.0 ± 2.0	pci/l 2.5	146.1 ± 22.7	pci/l 17.4
Gross Beta.....	3.3 ± 1.7	pci/l 2.5	8.7 ± 2.2	pci/l 2.8	134.3 ± 17.7	pci/l 17.9
Uranium 233, 234.....	0.39 ± 0.39	pci/l 0	3.05 ± 1.09	pci/l 0.43	84.9 ± 7.1	pci/l 0.69
Uranium 235.....	0.10 ± 0.19	pci/l 0	0.09 ± 0.32	pci/l 0.43	2.38 ± 1.26	pci/l 0.69
Uranium 238.....	0.10 ± 0.19	pci/l 0	1.85 ± 0.87	pci/l 0.43	49.7 ± 5.5	pci/l 0.69
Strontium 89, 90.....	0.11 ± 0.47	pci/l 0.76	0.68 ± 0.42	pci/l 0.61	0.49 ± 0.44	pci/l 0.67
Plutonium 239, 240.....	0.014 ± 0.008	pci/l 0	0.007 ± 0.008	pci/l 0.010	0.017 ± 0.011	pci/l 0
Americium 241.....	0.006 ± 0.012	pci/l 0	0.012 ± 0.014	pci/l 0	0.020 ± 0.010	pci/l 0
Cesium 137.....	0.12 ± 0.44	pci/l 0.71	-0.18 ± 0.38	pci/l 0.65	-0.11 ± 0.46	pci/l 0.77
Tritium.....	40 ± 300	pci/l 500	290 ± 280	pci/l 430	3430 ± 450	pci/l 530
Radium 226.....						
Radium 228.....						

Client: ROCKWELL (ROCKY FLATS)

Sample Information

	9003-803-6	9003-803-7
RFW Batch ID:		
Customer ID:	SH095007	SH093007
Collection Date:	10/10/89	10/10/89
Matrix:	Water	Water

Radio Chemistry

Gross Alpha.....	120.4 ± 21.2	pci/l 17.7	12.3 ± 2.7	pci/l 2.5
Gross Beta.....	141.1 ± 18.8	pci/l 20.2	13.4 ± 2.4	pci/l 2.7
Uranium 233, 234.....	70.2 ± 4.9	pci/l 0.40	4.22 ± 1.17	pci/l 0.36
Uranium 235.....	2.83 ± 1.02	pci/l 0.40	0.08 ± 0.27	pci/l 0.36
Uranium 238.....	45.3 ± 4.0	pci/l 0.40	5.31 ± 1.31	pci/l 0.36
Strontium 89, 90.....	0.26 ± 0.42	pci/l 0.65	0.77 ± 0.44	pci/l 0.65
Plutonium 239, 240.....	0.010 ± 0.008	pci/l 0.009	0.007 ± 0.005	pci/l 0
Americium 241.....			0.010 ± 0.008	pci/l 0
Cesium 137.....	0.07 ± 0.42	pci/l 0.68	0.07 ± 0.41	pci/l 0.67
Tritium.....	2730 ± 400	pci/l 500	-40 ± 250	pci/l 430
Radium 226.....				
Radium 228.....				

000 012